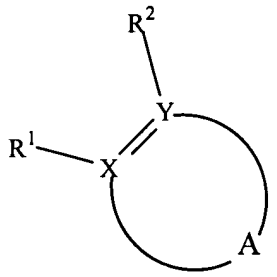
This listing of claims will replace all prior versions, and listings, of claims in the application.

1. (Original) A compound of Formula (I)



(I)

or a pharmaceutically acceptable salt or solvate thereof

wherein

R^1 is H, C_{1-6}alkyl, C_{1-6}haloalkyl, C_{1-6}alkoxy, C_{1-6}thioalkyl, cyano, halo, C_{3-7}cycloalkyl, -C_{1-6}alkylene-C_{3-7}cycloalkyl, C_{2-6}alkenyl or C_{3-6}alkynyl;

R^2 is $C(D)NR^3R^4$, D'- D''$(R^3)(R^4)$ or $CH_2N\ R^3R^4$

D' is CH_2 or a bond;

D'' is C, C-OH or CH

wherein

said C is attached to R^3 by a single or double bond;

said C is attached to R^4 by a single or double bond;

provided that

C is not attached to both R^3 and R^4 by double bonds;

said CH is attached to R^3 and R^4 by single bonds;

said C of C-OH is attached to R^3 and R^4 by single bonds;

D is O or S;

R^3 and R^4 are each independently selected from the group consisting of H, C_{1-6}alkyl, C_{1-6}haloalkyl, -C_{1-6}hydroxyalkyl, -C_{1-4}alkylene-O-C_{1-4}alkyl, -C_{1-3}alkylene-C_{1-6}thioalkyl, -C_{2-6}alkylidene-$(C_{1-4}$alkoxy$)_2$, C_{3-7}cycloalkyl, -C_{1-6}alkylene-C_{3-7}cycloalkyl, C_{2-6}alkenyl, C_{3-6}alkynyl, -C_{1-6}alkylene-CN, -C_{1-6}alkylene-heterocyclo and -C_{1-6}alkylene-aryl;

wherein said aryl of said -C_{1-6}alkylene-aryl is optionally substituted with one to three of the same or different substituents selected from the group consisting of fluoro, chloro, bromo, cyano, nitro, C_{1-4}alkyl and C_{1-3}alkoxy;

or

R^3 and R^4 together with the nitrogen to which they are attached form a five or six-membered heterocycle,

said heterocycle optionally containing one additional heteroatom selected from the group consisting of N, S and O; and

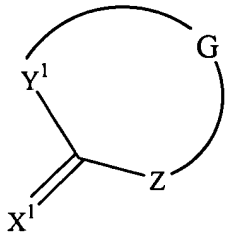
said heterocycle optionally substituted with one or more groups selected from the group consisting of C_{1-6}alkyl, C_{1-6}alkoxy, aryl, -C_{1-4}alkylene-aryl, pyridyl and halogen;

wherein said aryl of said -C_{1-4}alkylene-aryl is optionally substituted with one to three of the same or different substituents selected from the group consisting of fluoro, chloro, bromo, cyano, nitro and C_{1-3}alkoxy;

X is C;

Y is C;

A is

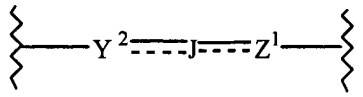


wherein

X^1 is N and is attached to X;

Y^1 is N and is attached to Y;

G is



wherein

Y^2 is CE^1 and is attached to Y^1;

J is a bond ;

Z^1 is CE^3 and is attached to Z;

wherein

E^1 and E^3 together form $N(CH)_3$,

optionally substituted with halogen, -CN, C_1-C_4alkyl, C_3-C_6cycloalkyl, substituted or unsubstituted phenyl, hydroxy, C_1-C_4alkoxy, SH, C_1-C_4thioalkyl, NH_2, $NH(C_1$-C_4alkyl) or $N(C_1$-C_4alkyl)$_2$;

Z is N-V, wherein V is phenyl, 2-pyridyl or 3-pyridyl substituted with two to three of the same or different substitutents selected from the group consisting of C_{1-4}alkyl, C_{1-4}alkoxy, C_{1-6}thioalkyl, C_{1-4}haloalkyl, halogen, $N(C_1$-C_4alkyl)$_2$ and CN.

2. (Original) A compound according to claim 1 wherein V is phenyl or 3-pyridyl and is substituted with two to three of the same or different substitutents selected from the group consisting of C_{1-4}alkyl, C_{1-4}alkoxy, C_{1-6}thioalkyl, C_{1-4}haloalkyl, halogen, $N(C_1$-C_4alkyl)$_2$ and CN; said substituents attached at the 2, 4 or 6-positions of said phenyl or said 3-pyridyl.

3. (Original) A compound according to claim 1 wherein V is 2-pyridyl and is substituted with two of the same or different substitutents selected from the group consisting of C_{1-4}alkyl, C_{1-4}alkoxy, C_{1-6}thioalkyl, C_{1-4}haloalkyl, halogen, $N(C_1-C_4alkyl)_2$ and CN; said substituents attached at the 3 and 5-positions of said 2-pyridyl.

4. (Original) A compound according to claim 1 wherein R^1 is C_{1-6}alkyl or C_{1-6}haloalkyl.

5. (Original) A compound according to claim 1 wherein R^1 is methyl or trifluoromethyl.

6. (Original) A compound according to claim 1 wherein R^2 is $C(D)NR^3R^4$ and D is O.

7. (Original) A compound according to claim 1 wherein R^2 is $CH_2N\ R^3R^4$.

8. (Original) A compound according to claim 1 wherein R^2 is D'- D''$(R^3)(R^4)$, D is a bond and D'' is C-OH.

9. (Original) A compound according to claim 1 wherein R^2 is D'- D''$(R^3)(R^4)$, D is a bond and D'' is C or CH.

10. (Original) A compound according to claim 1 wherein R^3 and R^4 are each independently selected from the group consisting of H, C_{1-6}alkyl, C_{1-6}haloalkyl, -C_{1-6}hydroxyalkyl, -C_{1-4}alkylene-O-C_{1-4}alkyl, -C_{1-3}alkylene-C_{1-6}thioalkyl, -C_{2-6}alkylidene-$(C_{1-4}alkoxy)_2$, C_{3-7}cycloalkyl, -C_{1-6}alkylene-C_{3-7}cycloalkyl, C_{2-6}alkenyl, C_{3-6}alkynyl and -C_{1-6}alkylene-CN.

11. (Original) A compound according to claim 1 wherein R^3 and R^4 together with the nitrogen to which they are attached form a five or six-membered heterocycle.

12. (Original) A compound according to claim 1 wherein V is 2, 4, 6-trimethylphenyl.

13. (Original) A compound according to claim 1 wherein V is 2,4-dichlorophenyl.

14. (Original) A compound according to claim 1 wherein E^1 and E^3 together form $N(CH)_3$ optionally substituted with halogen, methoxy, methyl or nitrile.

15. (Original) A compound according to claim 1 wherein R^2 is $CH_2NR^3R^4$, R^3 is ethyl or propyl, R^4 is $-(CH_2)_2$-phenyl, E^1 and E^3 together form $N(CH)_3$ optionally substituted with halogen, methoxy, methyl or nitrile.

16. (Original) A compound according to claim 1 wherein R^2 is $CH_2NR^3R^4$, R^3 is ethyl or propyl, R^4 is $-(CH_2)_2$-phenyl, E^1 and E^3 together form $N(CH)_3$ optionally substituted with halogen.

17. (Original) A compound or pharmaceutically acceptable salt of solvate thereof selected from the group consisting of

Ethyl-[2-methyl-8-(2,4,6-trimethyl-phenyl)-8*H*-1,3a,7,8-tetraaza-cyclopenta[α]inden-3-ylmethyl]-phenethyl-amine;

Cyclobutylmethyl-[2-methyl-8-(2,4,6-trimethyl-phenyl)-8*H*-1,3a,7,8-tetraaza-cyclopenta[α]inden-3-ylmethyl]-propyl-amine;

[8-(2-Chloro-4,6-dimethyl-phenyl)-2-methyl-8H-1,3a,7,8-tetraaza-cyclopenta[a]inden-3-ylmethyl]-phenethyl-propyl-amine;

[8-(2-Chloro-4,6-dimethyl-phenyl)-2-methyl-8*H*-1,3a,7,8-tetraaza-cyclopenta[α]inden-3-ylmethyl]-cyclobutylmethyl-propyl-amine;

[8-(2-Chloro-4,6-dimethyl-phenyl)-2-methyl-8H-1,3a,7,8-tetraaza-cyclopenta[a]inden-3-ylmethyl]-ethyl-phenethyl-amine;

8-(2-Chloro-4,6-dimethyl-phenyl)-2-methyl-3-(3-phenyl-pyrrolidin-1-ylmethyl)-8*H*-1,3a,7,8-tetraaza-cyclopenta[α]indene;

Cyclopropylmethyl-propyl-[2-trifluoromethyl-8-(2,4,6-trimethyl-phenyl)-8H-1,3a,7,8-tetraaza-cyclopenta[a]inden-3-ylmethyl]-amine;

Phenethyl-[2-trifluoromethyl-8-(2,4,6-trimethyl-phenyl)-8H-1,3a,7,8-tetraaza-cyclopenta[a]inden-3-ylmethyl]-(3,3,3-trifluoro-propyl)-amine;

[8-(2-Chloro-4,6-dimethyl-phenyl)-2-methyl-8H-1,3a,7,8-tetraaza-cyclopenta[a]inden-3-ylmethyl]-phenethyl-propyl-amine;

[8-(2-Chloro-4,6-dimethyl-phenyl)-2-methyl-8H-1,3a,7,8-tetraaza-cyclopenta[a]inden-3-

ylmethyl]-ethyl-phenethyl-amine; and

[8-(2-Chloro-4,6-dimethyl-phenyl)-2-methyl-8*H*-1,3a,7,8-tetraaza-cyclopenta[*a*]inden-3-ylmethyl]-(2-methoxy-1-methoxymethyl-ethyl)-amine.

18. (Cancelled)